

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121418\
 Data File : BG038596.D
 Acq On : 15 Dec 2018 9:35
 Operator : JU/SJ
 Sample : PB115627BL
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115627BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG121218.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.144	345	350	355	rVB2	15635	23970	1.54%	0.329%
2	5.608	422	429	438	rBV	330761	548461	35.22%	7.523%
3	7.212	694	702	712	rBV	327043	590385	37.91%	8.099%
4	7.589	758	766	777	rBV	309291	562510	36.12%	7.716%
5	8.059	839	846	855	rVB	77270	132672	8.52%	1.820%
6	8.382	894	901	910	rVB	246932	438300	28.15%	6.012%
7	9.234	1039	1046	1055	rBV	196137	369524	23.73%	5.069%
8	10.879	1318	1326	1335	rBV	89443	176348	11.32%	2.419%
9	13.317	1733	1741	1751	rBV	513373	857702	55.08%	11.765%
10	14.698	1969	1976	1985	rVB	152185	253714	16.29%	3.480%
11	16.184	2223	2229	2246	rBV	374393	621218	39.89%	8.522%
12	17.442	2436	2443	2456	rBV2	212333	345725	22.20%	4.742%
13	20.044	2880	2886	2894	rBV	1175698	1557248	100.00%	21.361%
14	21.719	3164	3171	3184	rVB	254134	427309	27.44%	5.862%
15	25.015	3721	3732	3743	rVB2	125783	384901	24.72%	5.280%

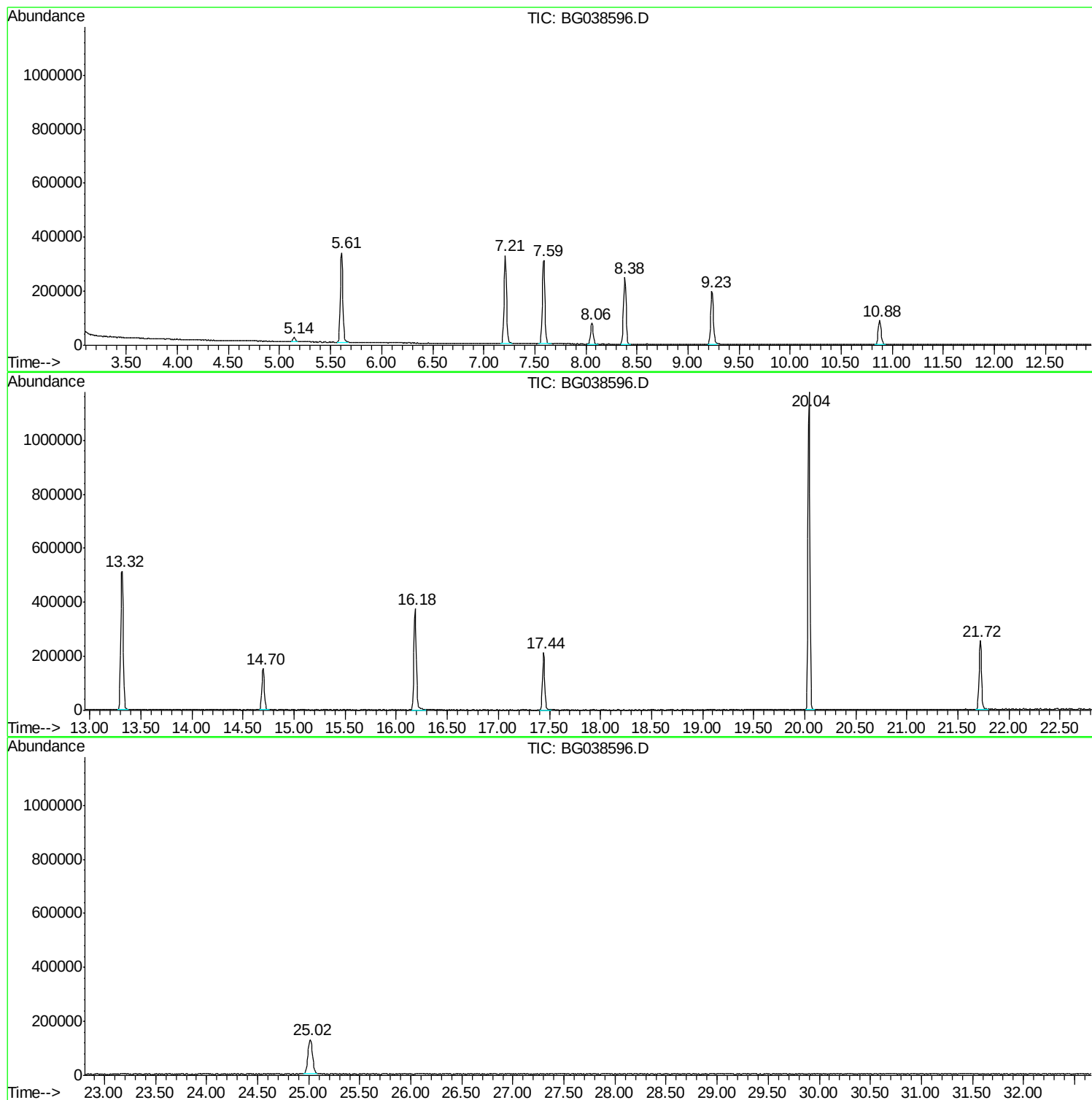
Sum of corrected areas: 7289987

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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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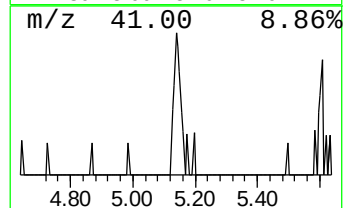
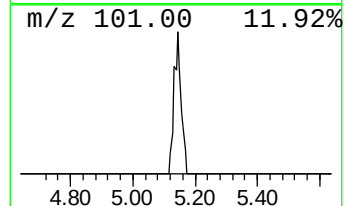
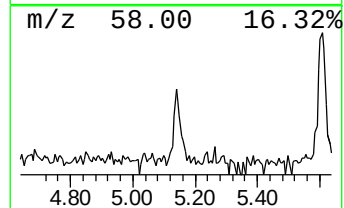
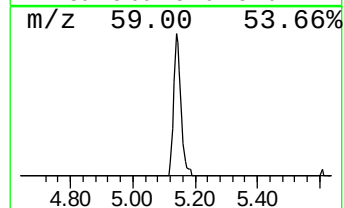
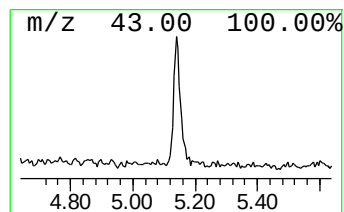
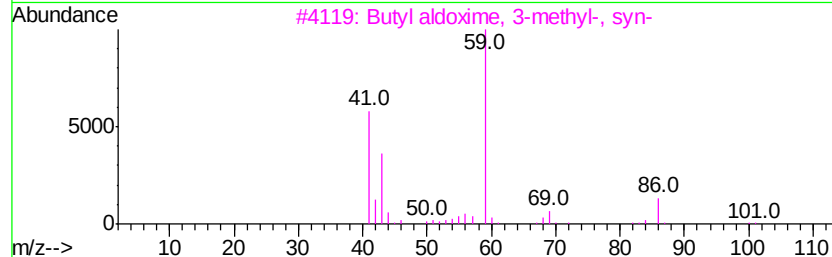
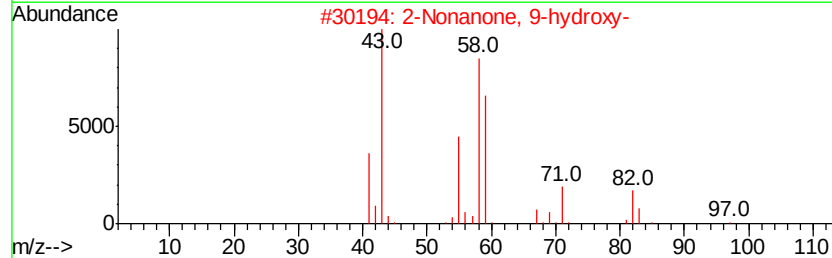
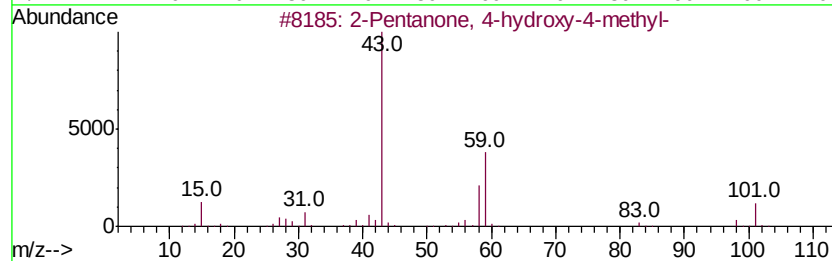
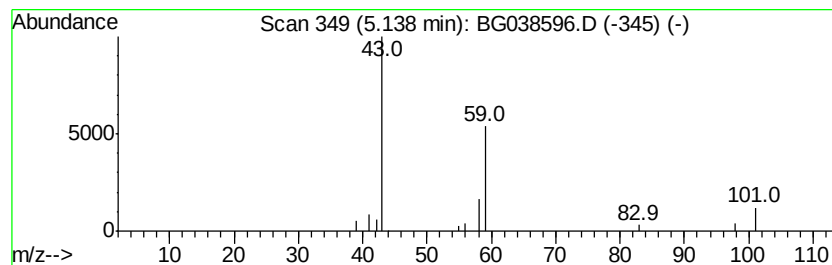
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	3.61 ng	23970	1,4-Dichlorobenzene-d4	8.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2		000123-42-2	50
2	2-Nonanone, 9-hydroxy-	158	C9H18O2		025368-56-3	36
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO		005780-40-5	23
4	2,3-Butanedione, monooxime	101	C4H7NO2		000057-71-6	9
5	1-Propen-2-ol, acetate	100	C5H8O2		000108-22-5	9



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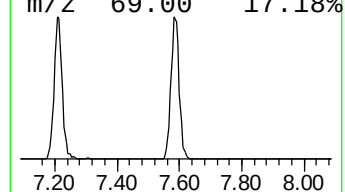
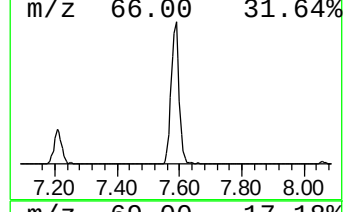
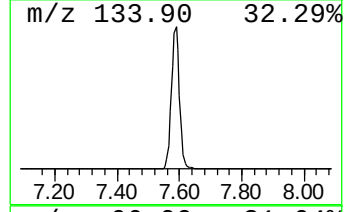
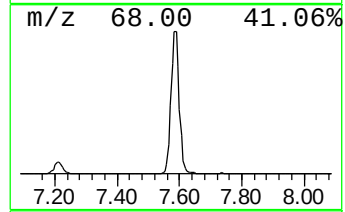
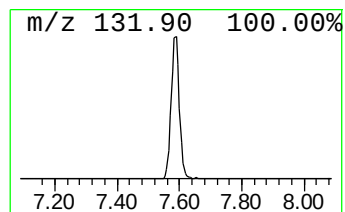
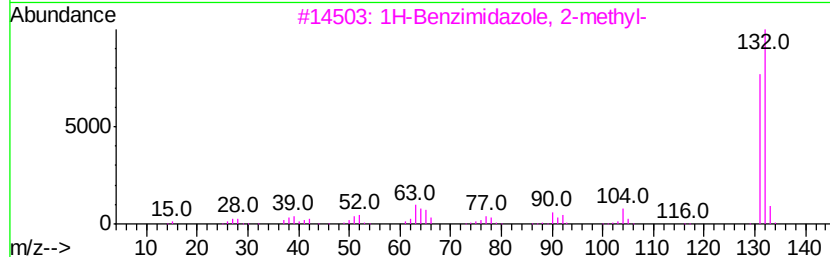
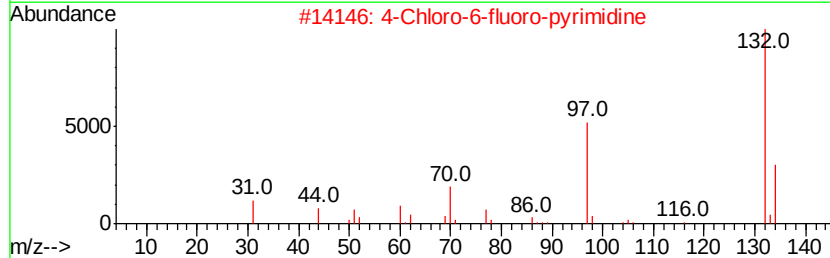
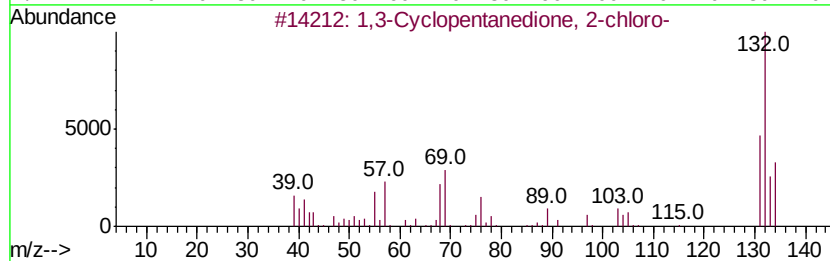
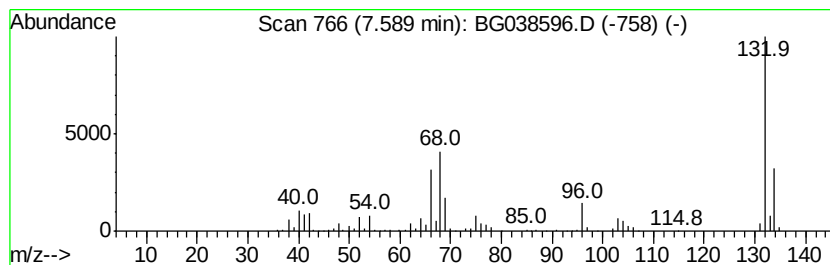
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Peak Number 2 unknown7.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	84.80 ng	562510	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	37
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	3.6	ng	23970	1	8.06	132672	20.0
unknown7.59	7.59	84.8	ng	562510	1	8.06	132672	20.0